Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF072517\
Data File : BF097024.D
Acq On : 25 Jul 2017 12:58
Operator : SJ/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF072517

Quant Time: Jul 25 14:12:00 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 25 14:11:51 2017
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	93	0.00
2		1,4-Dioxane	40.000	41.794	-4.5	106	0.00
3		Pyridine	40.000	41.872	-4.7	89	0.00
4		n-Nitrosodimethylamine	40.000	40.449	-1.1	92	0.00
5	S	2-Fluorophenol	80.000	77.687	2.9	94	0.00
6		Aniline	40.000	39.728	0.7	91	0.00
7	S	Phenol-d6	80.000	79.034	1.2	92	0.00
8		2-Chlorophenol	40.000	40.128	-0.3	94	0.00
9		Benzaldehyde	40.000	39.117	2.2	94	0.00
10	C	Phenol	40.000	41.059	-2.6	93	0.00
11		bis(2-Chloroethyl)ether	40.000	41.723	-4.3	96	0.00
12		1,3-Dichlorobenzene	40.000	40.107	-0.3	94	0.00
13	C	1,4-Dichlorobenzene	40.000	40.425	-1.1	100	0.00
14		1,2-Dichlorobenzene	40.000	41.282	-3.2	101	0.00
15		Benzyl Alcohol	40.000	39.885	0.3	101	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.699	0.8	99	0.00
17		2-Methylphenol	40.000	39.202	2.0	97	0.00
18		Hexachloroethane	40.000	38.423	3.9	93	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.291	1.8	98	0.00
20		3+4-Methylphenols	40.000	40.589	-1.5	101	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	96	0.00
22		Acetophenone	40.000	39.871	0.3	98	0.00
23	S	Nitrobenzene-d5	80.000	76.437	4.5	95	0.00
24		Nitrobenzene	40.000	39.114	2.2	94	0.00
25		Isophorone	40.000	36.261	9.3	91	0.00
26	C	2-Nitrophenol	40.000	40.726	-1.8	97	0.00
27		2,4-Dimethylphenol	40.000	37.445	6.4	86	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.993	2.5	93	0.00
29	C	2,4-Dichlorophenol	40.000	40.628	-1.6	94	0.00
30		1,2,4-Trichlorobenzene	40.000	40.590	-1.5	99	0.00
31		Naphthalene	40.000	39.740	0.6	100	0.00
32		Benzoic acid	40.000	40.578	-1.4	92	0.00
33		4-Chloroaniline	40.000	36.943	7.6	87	0.00
34	C	Hexachlorobutadiene	40.000	37.523	6.2	91	0.00
35		Caprolactam	40.000	36.426	8.9	84	0.00
36	C	4-Chloro-3-methylphenol	40.000	36.870	7.8	86	0.00
37		2-Methylnaphthalene	40.000	36.685	8.3	88	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	89	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	40.696	-1.7	89	0.00
40	P	Hexachlorocyclopentadiene	40.000	40.057	-0.1	87	0.00
41	S	2,4,6-Tribromophenol	80.000	79.804	0.2	92	0.00
42	C	2,4,6-Trichlorophenol	40.000	38.938	2.7	84	0.00
43		2,4,5-Trichlorophenol	40.000	40.159	-0.4	86	0.00
44	S	2-Fluorobiphenyl	80.000	82.182	-2.7	91	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF072517\
Data File : BF097024.D
Acq On : 25 Jul 2017 12:58
Operator : SJ/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF072517

Quant Time: Jul 25 14:12:00 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 25 14:11:51 2017
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	40.430	-1.1	89	0.00
46		2-Chloronaphthalene	40.000	39.650	0.9	87	0.00
47		2-Nitroaniline	40.000	38.086	4.8	78	0.00
48		Acenaphthylene	40.000	39.188	2.0	93	0.00
49		Dimethylphthalate	40.000	39.285	1.8	93	0.00
50		2,6-Dinitrotoluene	40.000	40.924	-2.3	94	0.00
51	C	Acenaphthene	40.000	39.070	2.3	91	0.00
52		3-Nitroaniline	40.000	38.532	3.7	90	0.00
53	P	2,4-Dinitrophenol	40.000	42.200	-5.5	90	0.00
54		Dibenzofuran	40.000	39.621	0.9	92	0.00
55	P	4-Nitrophenol	40.000	41.816	-4.5	89	0.00
56		2,4-Dinitrotoluene	40.000	39.677	0.8	92	0.00
57		Fluorene	40.000	40.706	-1.8	93	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	40.712	-1.8	93	0.00
59		Diethylphthalate	40.000	40.686	-1.7	96	0.00
60		4-Chlorophenyl-phenylether	40.000	40.864	-2.2	93	0.00
61		4-Nitroaniline	40.000	40.873	-2.2	91	0.00
62		Azobenzene	40.000	39.486	1.3	84	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	91	0.00
64		4,6-Dinitro-2-methylphenol	40.000	42.441	-6.1	95	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.144	7.1	94	0.00
66		4-Bromophenyl-phenylether	40.000	40.358	-0.9	100	0.00
67		Hexachlorobenzene	40.000	39.075	2.3	97	0.00
68		Atrazine	40.000	40.124	-0.3	102	0.00
69	C	Pentachlorophenol	40.000	38.843	2.9	87	0.00
70		Phenanthrene	40.000	39.758	0.6	97	0.00
71		Anthracene	40.000	38.499	3.8	94	0.00
72		Carbazole	40.000	38.675	3.3	97	0.00
73		Di-n-butylphthalate	40.000	39.271	1.8	97	0.00
74	C	Fluoranthene	40.000	38.157	4.6	98	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	91	0.00
76		Benzidine	40.000	41.504	-3.8	88	0.00
77		Pyrene	40.000	43.182	-8.0	105	0.00
78	S	Terphenyl-d14	80.000	82.506	-3.1	101	0.00
79		Butylbenzylphthalate	40.000	45.092	-12.7	105	0.00
80		Benzo(a)anthracene	40.000	39.882	0.3	95	0.00
81		3,3'-Dichlorobenzidine	40.000	38.918	2.7	93	0.00
82		Chrysene	40.000	39.718	0.7	95	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	40.960	-2.4	97	0.00
84	c	Di-n-octyl phthalate	40.000	41.327	-3.3	96	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	36.207	9.5	91	0.00
86	I	Perylene-d12	20.000	20.000	0.0	91	0.00
87		Benzo(b)fluoranthene	40.000	38.407	4.0	91	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF072517\
Data File : BF097024.D
Acq On : 25 Jul 2017 12:58
Operator : SJ/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF072517

Quant Time: Jul 25 14:12:00 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 25 14:11:51 2017
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	39.096	2.3	94	0.00
89 C	Benzo(a)pyrene	40.000	39.683	0.8	96	0.00
90	Dibenzo(a,h)anthracene	40.000	36.570	8.6	91	0.00
91	Benzo(g,h,i)perylene	40.000	36.766	8.1	91	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 0